



# Full wwPDB X-ray Structure Validation Report ⓘ

May 29, 2020 – 02:35 am BST

PDB ID : 3G8U  
Title : DNA binding domain:GilZ 16bp complex-5  
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Deposited on : 2009-02-12  
Resolution : 1.90 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Xtriage (Phenix) : 1.13  
EDS : 2.11  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.11

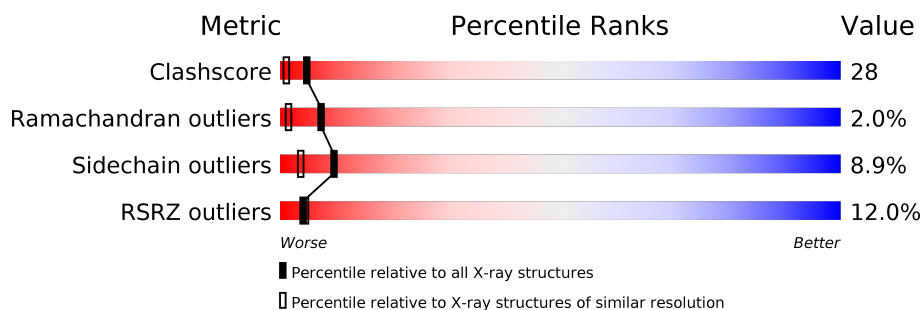
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.90 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)
RSRZ outliers	127900	6082 (1.90-1.90)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	90	<div> <div>11%</div> <div> <div></div> <div>52%</div> <div>30%</div> <div>6%</div> <div>12%</div> </div> </div>
1	B	90	<div> <div>12%</div> <div> <div></div> <div>56%</div> <div>19%</div> <div>• •</div> <div>20%</div> </div> </div>
2	D	16	<div> <div>6%</div> <div> <div></div> <div>63%</div> <div>38%</div> </div> </div>
3	C	16	<div> <div>25%</div> <div> <div></div> <div>38%</div> <div>38%</div> </div> </div>

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 1914 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Glucocorticoid receptor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	79	Total	C	N	O	S	0	3	0
			627	379	125	110	13			
1	B	72	Total	C	N	O	S	0	0	0
			549	335	105	97	12			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	436	GLY	-	EXPRESSION TAG	UNP P06536
A	437	SER	-	EXPRESSION TAG	UNP P06536
A	438	HIS	-	EXPRESSION TAG	UNP P06536
A	439	MET	-	EXPRESSION TAG	UNP P06536
B	436	GLY	-	EXPRESSION TAG	UNP P06536
B	437	SER	-	EXPRESSION TAG	UNP P06536
B	438	HIS	-	EXPRESSION TAG	UNP P06536
B	439	MET	-	EXPRESSION TAG	UNP P06536

- Molecule 2 is a DNA chain called DNA (5'-D(\*TP\*GP\*GP\*AP\*AP\*CP\*CP\*CP\*AP\*AP\*TP\*GP\*TP\*TP\*CP\*T)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	16	Total	C	N	O	P	0	0	0
			323	156	57	95	15			

- Molecule 3 is a DNA chain called DNA (5'-D(\*AP\*AP\*GP\*AP\*AP\*CP\*AP\*TP\*TP\*GP\*GP\*GP\*TP\*TP\*CP\*C)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	16	Total	C	N	O	P	0	0	0
			327	157	62	93	15			

- Molecule 4 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	2	Total 2	Zn 2	0	0
4	A	2	Total 2	Zn 2	0	0

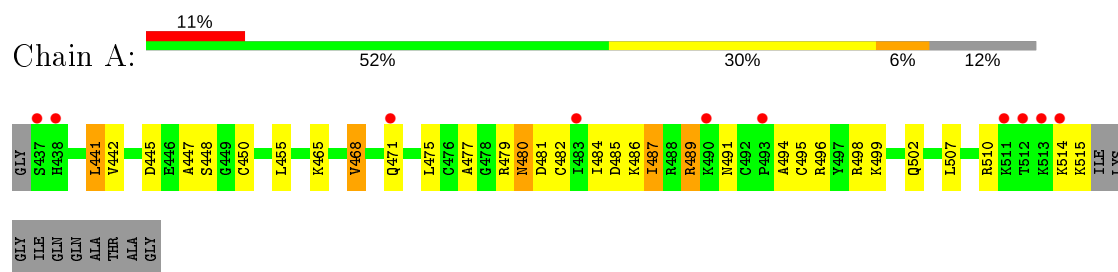
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	45	Total 45	O 45	0	0
5	B	25	Total 25	O 25	0	0
5	D	12	Total 12	O 12	0	0
5	C	2	Total 2	O 2	0	0

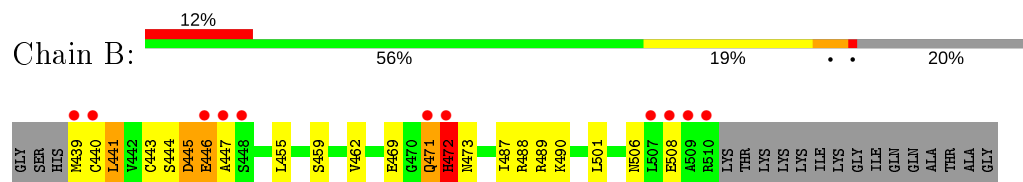
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

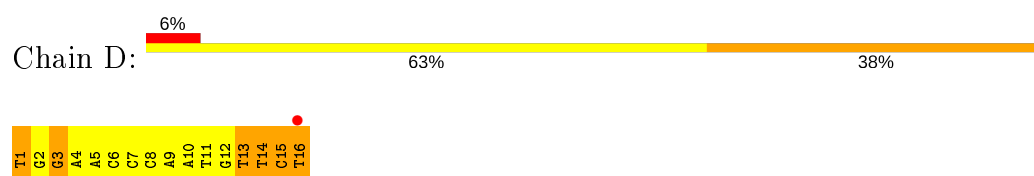
#### • Molecule 1: Glucocorticoid receptor



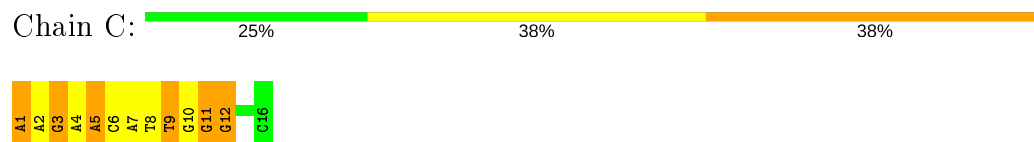
#### • Molecule 1: Glucocorticoid receptor



#### • Molecule 2: DNA (5'-D(\*TP\*GP\*GP\*AP\*AP\*CP\*CP\*CP\*AP\*AP\*TP\*GP\*TP\*TP\*CP\*T)-3')



#### • Molecule 3: DNA (5'-D(\*AP\*AP\*GP\*AP\*AP\*CP\*AP\*TP\*TP\*GP\*GP\*GP\*TP\*TP\*CP\*C)-3')



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	116.94Å 38.60Å 98.52Å 90.00° 123.51° 90.00°	Depositor
Resolution (Å)	29.20 – 1.90 29.20 – 1.90	Depositor EDS
% Data completeness (in resolution range)	86.8 (29.20-1.90) 86.8 (29.20-1.90)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	0.07	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.15 (at 1.91Å)	Xtriage
Refinement program	PHENIX (phenix.refine)	Depositor
R, $R_{free}$	0.210 , 0.236 0.205 , (Not available)	Depositor DCC
$R_{free}$ test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	41.5	Xtriage
Anisotropy	0.449	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 74.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.004 for h,-k,-h-l	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	1914	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	82.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.71% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: ZN

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 5$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z  > 5$	RMSZ	$\# Z  > 5$
1	A	0.99	3/635 (0.5%)	0.90	2/847 (0.2%)
1	B	0.61	0/556	0.71	0/742
2	D	1.26	3/361 (0.8%)	2.08	19/555 (3.4%)
3	C	1.02	0/367	2.17	23/565 (4.1%)
All	All	0.96	6/1919 (0.3%)	1.50	44/2709 (1.6%)

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	12	DG	C5'-C4'	7.73	1.59	1.51
2	D	12	DG	P-O5'	7.04	1.66	1.59
1	A	495	CYS	CB-SG	6.34	1.93	1.82
2	D	11	DT	C1'-N1	6.33	1.57	1.49
1	A	494	ALA	CA-CB	5.35	1.63	1.52
1	A	482	CYS	CB-SG	5.13	1.91	1.82

All (44) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	14	DT	O4'-C1'-N1	-12.26	99.42	108.00
3	C	1	DA	C2-N3-C4	-10.36	105.42	110.60
3	C	1	DA	C5-C6-N1	-9.15	113.12	117.70
3	C	5	DA	O4'-C1'-N9	8.66	114.06	108.00
3	C	3	DG	O4'-C1'-N9	8.64	114.05	108.00
3	C	9	DT	P-O3'-C3'	8.16	129.50	119.70
2	D	16	DT	O4'-C4'-C3'	-8.00	101.20	106.00
3	C	1	DA	C1'-O4'-C4'	-7.81	102.29	110.10
2	D	12	DG	O4'-C4'-C3'	-7.43	101.53	104.50
2	D	16	DT	C1'-O4'-C4'	-7.42	102.68	110.10
3	C	2	DA	O5'-P-OP2	-7.13	99.29	105.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	12	DG	O4'-C1'-N9	7.07	112.94	108.00
2	D	16	DT	O4'-C1'-N1	7.06	112.94	108.00
3	C	3	DG	C3'-C2'-C1'	-6.65	94.52	102.50
3	C	1	DA	O4'-C1'-C2'	6.52	111.12	105.90
2	D	14	DT	N3-C4-O4	6.43	123.76	119.90
2	D	11	DT	N3-C4-O4	6.41	123.74	119.90
3	C	11	DG	OP1-P-O3'	6.27	119.00	105.20
1	A	496	ARG	NE-CZ-NH1	-6.19	117.21	120.30
2	D	13	DT	O4'-C1'-N1	5.97	112.18	108.00
2	D	16	DT	C4'-C3'-C2'	-5.81	97.87	103.10
3	C	5	DA	C3'-C2'-C1'	-5.78	95.56	102.50
1	A	489	ARG	NE-CZ-NH2	-5.78	117.41	120.30
2	D	3	DG	C1'-O4'-C4'	-5.70	104.40	110.10
3	C	5	DA	C4'-C3'-C2'	-5.65	98.01	103.10
3	C	8	DT	O4'-C1'-N1	5.62	111.93	108.00
2	D	1	DT	C5-C4-O4	-5.58	120.99	124.90
2	D	15	DC	O4'-C1'-C2'	-5.47	101.52	105.90
3	C	5	DA	P-O3'-C3'	5.35	126.12	119.70
3	C	11	DG	O4'-C1'-C2'	5.35	110.18	105.90
2	D	3	DG	O4'-C1'-C2'	-5.35	101.62	105.90
3	C	1	DA	N1-C2-N3	5.33	131.96	129.30
2	D	11	DT	C5-C4-O4	-5.32	121.18	124.90
2	D	11	DT	N1-C1'-C2'	5.30	122.68	112.60
3	C	1	DA	O4'-C1'-N9	5.29	111.70	108.00
3	C	1	DA	C6-N1-C2	5.27	121.77	118.60
3	C	1	DA	C5-C6-N6	5.26	127.91	123.70
2	D	1	DT	O4'-C1'-N1	5.24	111.67	108.00
2	D	11	DT	C5'-C4'-C3'	5.22	123.50	114.10
3	C	11	DG	P-O3'-C3'	-5.21	113.45	119.70
3	C	11	DG	P-O5'-C5'	-5.19	112.60	120.90
3	C	7	DA	N1-C6-N6	5.18	121.71	118.60
2	D	16	DT	C5'-C4'-C3'	5.07	123.23	114.10
3	C	12	DG	O4'-C1'-N9	5.03	111.52	108.00

There are no chirality outliers.

There are no planarity outliers.

## 5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen



atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	627	0	604	33	0
1	B	549	0	534	31	0
2	D	323	0	183	19	0
3	C	327	0	182	26	0
4	A	2	0	0	0	0
4	B	2	0	0	0	0
5	A	45	0	0	6	0
5	B	25	0	0	3	0
5	C	2	0	0	0	0
5	D	12	0	0	0	0
All	All	1914	0	1503	93	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 28.

All (93) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:1:DA:H8	3:C:1:DA:HO5'	1.14	0.96
2:D:15:DC:H3'	2:D:16:DT:H4'	1.48	0.94
3:C:4:DA:H2''	3:C:5:DA:H5'	1.59	0.84
1:A:450[B]:CYS:O	3:C:1:DA:H5'	1.87	0.74
2:D:7:DC:H42	3:C:11:DG:H1	1.38	0.70
3:C:4:DA:C2'	3:C:5:DA:H5'	2.23	0.69
1:A:487:ILE:O	1:A:487:ILE:HD13	1.94	0.68
2:D:15:DC:H2''	2:D:16:DT:H5'	1.77	0.67
1:B:447:ALA:CB	1:B:455:LEU:HD11	2.25	0.66
2:D:7:DC:N3	3:C:11:DG:N2	2.38	0.65
2:D:6:DC:H2''	2:D:7:DC:H5'	1.79	0.65
1:B:443:CYS:SG	1:B:445:ASP:HB2	2.37	0.65
2:D:5:DA:H2''	2:D:6:DC:O5'	1.99	0.63
1:B:439:MET:HA	1:B:446:GLU:HG2	1.80	0.62
3:C:1:DA:C8	3:C:1:DA:C5'	2.84	0.61
1:A:485:ASP:N	5:A:82:HOH:O	2.33	0.60
2:D:15:DC:C2'	2:D:16:DT:H5'	2.30	0.60
1:A:475:LEU:CD1	1:B:487:ILE:HG21	2.31	0.60
1:B:489:ARG:NH2	3:C:12:DG:H5'	2.17	0.60
2:D:9:DA:H2''	2:D:10:DA:OP2	2.02	0.60
1:A:484:ILE:C	5:A:82:HOH:O	2.40	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:472:HIS:HB2	5:B:52:HOH:O	2.01	0.59
1:A:485:ASP:HB3	5:A:82:HOH:O	2.03	0.59
1:A:447:ALA:CB	1:A:455:LEU:CD2	2.80	0.58
1:A:484:ILE:HG23	1:A:489:ARG:HA	1.85	0.58
1:A:475:LEU:HD12	1:B:487:ILE:HG21	1.84	0.58
1:B:506:ASN:OD1	1:B:508:GLU:HG2	2.04	0.58
1:A:487:ILE:HD13	1:A:487:ILE:C	2.25	0.57
1:B:489:ARG:NH2	3:C:12:DG:C5'	2.69	0.56
1:A:498:ARG:O	1:A:502:GLN:HG3	2.06	0.56
2:D:6:DC:H2'	2:D:7:DC:C6	2.41	0.56
3:C:1:DA:C8	3:C:1:DA:H5''	2.42	0.55
1:B:443:CYS:SG	1:B:489:ARG:HD2	2.47	0.55
1:B:489:ARG:HH21	3:C:12:DG:H5'	1.72	0.55
1:B:472:HIS:CB	5:B:52:HOH:O	2.55	0.54
1:A:475:LEU:HD12	1:B:487:ILE:CG2	2.38	0.54
1:B:473:ASN:N	5:B:52:HOH:O	2.41	0.53
1:A:510:ARG:HB2	5:A:34:HOH:O	2.08	0.53
2:D:1:DT:H2''	2:D:2:DG:C8	2.43	0.53
1:A:447:ALA:HB1	1:A:455:LEU:CD2	2.38	0.53
3:C:4:DA:H1'	3:C:5:DA:H5'	1.90	0.52
3:C:1:DA:C8	3:C:1:DA:O5'	2.62	0.51
2:D:16:DT:OP2	2:D:16:DT:C4'	2.57	0.51
1:B:447:ALA:CB	1:B:455:LEU:CD1	2.89	0.51
2:D:3:DG:H2''	2:D:4:DA:C8	2.46	0.51
1:A:445:ASP:OD2	1:A:486:LYS:HE3	2.11	0.50
1:B:440:CYS:O	1:B:444:SER:HA	2.11	0.50
1:B:489:ARG:HH21	3:C:12:DG:C5'	2.25	0.50
1:B:447:ALA:HB3	1:B:455:LEU:HD11	1.91	0.50
1:A:479[B]:ARG:O	1:A:479[B]:ARG:CG	2.59	0.50
2:D:8:DC:H2''	2:D:9:DA:C8	2.47	0.49
2:D:16:DT:OP2	2:D:16:DT:H4'	2.12	0.49
1:B:439:MET:CA	1:B:446:GLU:HG2	2.42	0.48
1:B:471:GLN:HG2	1:B:472:HIS:N	2.29	0.47
1:A:468:VAL:HG11	1:A:507:LEU:HD13	1.97	0.47
3:C:1:DA:H2'	3:C:1:DA:N3	2.31	0.46
3:C:5:DA:H1'	3:C:6:DC:C6	2.51	0.45
1:A:447:ALA:HB1	1:A:455:LEU:HD23	1.98	0.45
2:D:6:DC:H2''	2:D:7:DC:C5'	2.46	0.45
1:A:442:VAL:O	1:A:499:LYS:HE2	2.18	0.44
1:B:444:SER:C	1:B:445:ASP:O	2.55	0.44
1:A:514:LYS:O	1:A:515:LYS:CB	2.65	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:477:ALA:HA	1:B:488:ARG:HD2	2.00	0.44
1:A:450[A]:CYS:O	3:C:1:DA:H4'	2.18	0.44
1:B:472:HIS:N	1:B:472:HIS:ND1	2.66	0.43
3:C:9:DT:C2'	3:C:10:DG:C8	3.01	0.43
1:A:487:ILE:CD1	1:A:487:ILE:C	2.87	0.43
2:D:13:DT:C2'	2:D:14:DT:H72	2.49	0.43
3:C:4:DA:C1'	3:C:5:DA:H5'	2.48	0.43
2:D:13:DT:H2'	2:D:14:DT:H72	1.99	0.43
1:B:471:GLN:O	1:B:472:HIS:O	2.37	0.43
2:D:15:DC:C3'	2:D:16:DT:H4'	2.32	0.43
1:B:459:SER:O	1:B:462:VAL:HG22	2.18	0.43
1:B:471:GLN:O	1:B:472:HIS:C	2.57	0.43
1:A:479[B]:ARG:O	1:A:481:ASP:N	2.51	0.43
1:A:491:ASN:HD22	1:A:491:ASN:N	2.15	0.43
1:A:450[B]:CYS:O	3:C:1:DA:C5'	2.63	0.43
1:A:479[B]:ARG:HG2	1:A:479[B]:ARG:O	2.18	0.43
3:C:3:DG:H2'	3:C:3:DG:O5'	2.18	0.43
3:C:1:DA:H8	3:C:1:DA:O5'	1.88	0.42
1:B:490:LYS:HD2	3:C:11:DG:H4'	2.00	0.42
1:B:439:MET:N	1:B:446:GLU:HG2	2.35	0.42
1:A:479[B]:ARG:O	1:A:480:ASN:C	2.57	0.42
1:A:448:SER:HB3	5:A:40:HOH:O	2.20	0.42
2:D:7:DC:N4	3:C:11:DG:H1	2.10	0.42
1:B:489:ARG:NH2	3:C:12:DG:H5''	2.33	0.42
1:B:508:GLU:CG	1:B:508:GLU:O	2.68	0.42
1:A:485:ASP:CA	5:A:82:HOH:O	2.68	0.41
1:A:484:ILE:HG23	1:A:489:ARG:CB	2.50	0.41
1:B:441:LEU:CD2	1:B:441:LEU:N	2.84	0.40
3:C:4:DA:H2''	3:C:5:DA:OP2	2.21	0.40
1:A:441:LEU:HD12	1:A:441:LEU:HA	1.77	0.40
1:A:507:LEU:HA	1:A:507:LEU:HD12	1.69	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles ⓘ

### 5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	80/90 (89%)	75 (94%)	4 (5%)	1 (1%)	12	4
1	B	70/90 (78%)	62 (89%)	6 (9%)	2 (3%)	4	1
All	All	150/180 (83%)	137 (91%)	10 (7%)	3 (2%)	7	1

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	480	ASN
1	B	445	ASP
1	B	472	HIS

### 5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	67/73 (92%)	62 (92%)	5 (8%)	13	5
1	B	59/73 (81%)	53 (90%)	6 (10%)	7	2
All	All	126/146 (86%)	115 (91%)	11 (9%)	9	4

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	441	LEU
1	A	465	LYS
1	A	468	VAL
1	A	471	GLN
1	A	487	ILE
1	B	441	LEU
1	B	446	GLU
1	B	469	GLU
1	B	471	GLN
1	B	472	HIS

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Mol	Chain	Res	Type
1	B	501	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	451	HIS
1	A	471	GLN
1	A	491	ASN
1	A	506	ASN
1	B	480	ASN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	79/90 (87%)	0.77	10 (12%) 3 4	35, 55, 113, 172	0
1	B	72/90 (80%)	1.02	11 (15%) 2 2	36, 79, 169, 198	0
2	D	16/16 (100%)	0.39	1 (6%) 20 22	71, 103, 128, 132	0
3	C	16/16 (100%)	0.47	0 100 100	73, 97, 137, 143	0
All	All	183/212 (86%)	0.81	22 (12%) 4 4	35, 77, 136, 198	0

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	439	MET	6.1
1	B	509	ALA	4.8
1	B	508	GLU	3.8
1	B	507	LEU	3.8
1	B	447	ALA	3.8
1	B	510	ARG	3.8
1	A	437	SER	3.2
1	A	512	THR	3.2
2	D	16	DT	3.2
1	A	438	HIS	3.1
1	A	493	PRO	3.0
1	B	472	HIS	2.9
1	A	511	LYS	2.9
1	B	448	SER	2.8
1	A	514	LYS	2.6
1	B	446	GLU	2.5
1	B	471	GLN	2.5
1	A	513	LYS	2.4
1	A	483	ILE	2.4
1	A	490	LYS	2.4
1	A	471	GLN	2.4
1	B	440	CYS	2.2

## 6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

## 6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
4	ZN	B	526	1/1	0.96	0.09	88,88,88,88	0
4	ZN	A	526	1/1	0.99	0.12	57,57,57,57	0
4	ZN	A	527	1/1	1.00	0.13	38,38,38,38	0
4	ZN	B	527	1/1	1.00	0.14	43,43,43,43	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.